

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE101618\
 Data File : BE097923.D
 Acq On : 16 Oct 2018 23:09
 Operator : SJ/JU
 Sample : J5393-10
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 RE137-700FT-20181009

Quant Time: Oct 17 16:26:54 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE101618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 16 15:36:44 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	1932	0.40	ng	0.00
7) Naphthalene-d8	10.69	136	8085	0.40	ng	0.00
13) Acenaphthene-d10	14.56	164	5431	0.40	ng	0.00
19) Phenanthrene-d10	17.29	188	17868	0.40	ng	-0.01
27) Chrysene-d12	21.48	240	24343	0.40	ng	-0.01
34) Perylene-d12	24.03	264	19329	0.40	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.47	112	1159	0.19	ng	0.00
5) Phenol-d6	7.06	99	984	0.11	ng	0.00
8) Nitrobenzene-d5	9.04	82	2769	0.36	ng	0.00
11) 2-Methylnaphthalene-d10	12.28	152	5049	0.38	ng	-0.02
14) 2,4,6-Tribromophenol	16.04	330	1155	0.40	ng	0.00
15) 2-Fluorobiphenyl	13.17	172	8836	0.40	ng	0.00
25) Fluoranthene-d10	19.33	212	105696	0.44	ng	0.00
29) Terphenyl-d14	19.92	244	35040	0.63	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.38	88	18236	5.414	ng	95
6) bis(2-Chloroethyl)ether	7.30	93	451	0.059	ng	89

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE101618\
Data File : BE097923.D
Acq On : 16 Oct 2018 23:09
Operator : SJ/JU
Sample : J5393-10
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_E
ClientSampleId :
RE137-700FT-20181009

Quant Time: Oct 17 16:26:54 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE101618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 16 15:36:44 2018
Response via : Initial Calibration

